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Approximation Properties of Coarse Spaces by Algebraic Multigrid

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Abstract

This note provides an application of the “window”-based spectral AMG method (cf. [FVZ05] or [Va08]) for proving energy error estimates desired in providing coarse (upscaled) discretization of fairly general classes of PDEs. Computationally more efficient versions of the original window-based spectral AMG, as well as a new method utilizing local element matrices, are outlined summarizing the results from a forthcoming report.

1 The strong approximation property

We are given a s.p.d. $n \times n$ sparse matrix A and let $P : \mathbf{R}^{n_c} \mapsto \mathbf{R}^n$, $n_c < n$, be a given (rectangular) interpolation matrix.

We are interested in the following *strong approximation property*:

For any fine-grid vector $\mathbf{u} \in \mathbf{R}^n$ there is a coarse interpolant $P\mathbf{u}_c$ such that

$$\|A\| \|\mathbf{u} - P\mathbf{u}_c\|_A^2 \leq C_A \|A\mathbf{u}\|^2. \quad (1.1)$$

If the problem of our main interest

$$A\mathbf{u} = \mathbf{f},$$

comes from a finite element discretization of a PDE on a domain $\Omega \subset \mathbf{R}^d$ ($d = 2$ or 3), then $\mathbf{f} = (f_i)$ comes from a given r.h.s. function $f(\mathbf{x}) \in L_2(\Omega)$, where the entries f_i are computed as the following integral moments

$$f_i = (f, \varphi_i) \equiv \int_{\Omega} f(\mathbf{x}) \varphi_i d\mathbf{x}.$$

Above, φ_i runs over a basis of the fine-grid finite element space V_h associated with a triangulation of Ω with characteristic fine-grid mesh size h . For a nodal (Lagrangian)

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basis, the index “ i ” runs over the set of fine degrees of freedom $x_i \in \mathcal{N}_h$. The unknown \mathbf{u} stands for the coefficient vector of the finite element, Galerkin, approximation u_h to the solution of the underlined PDE posed variationally, i.e., $u_h \in V_h$ solves the discretized PDE in a variational form associated with a given bilinear form $a(\cdot, \cdot)$, stated as follows

$$a(u_h, \varphi) = (f, \varphi) \text{ for all } \varphi \in V_h.$$

As an example, we consider a second order self-adjoint elliptic bilinear form $a(u, \varphi) = \int_{\Omega} k(\mathbf{x}) \nabla u \cdot \nabla \varphi \, d\mathbf{x}$ for $u, \varphi \in H_0^1(\Omega)$ and a given positive coefficient function $k = k(\mathbf{x})$, $\mathbf{x} \in \Omega$, the given polygonal/polyhedral domain in \mathbf{R}^d , $d = 2$ or 3 . Using a standard piecewise linear conforming finite element space V_h on a quasiuniform triangulation \mathcal{T}_h , as it is well-known, the stiffness matrix $A = (a(\varphi_j, \varphi_i))$ computed from a nodal Lagrangian basis $\{\varphi_i\}_{\mathbf{x}_i \in \mathcal{N}_h}$ of V_h satisfies

$$\|A\| \simeq h^{d-2}. \quad (1.2)$$

The equivalence constants above generally depend on the variation $\frac{\max_{\mathbf{x} \in \Omega} k(\mathbf{x})}{\min_{\mathbf{x} \in \Omega} k(\mathbf{x})}$ but are mesh-independent.

Assume, that we have come up with a coarse space $V_H \subset V_h$ such that the coefficient vectors of functions in V_H viewed as elements of V_h can be represented as the range of an interpolation mapping P . We can define respective coarse basis functions by forming $P\mathbf{e}_{i_c}$ for each coarse coordinate vector $\mathbf{e}_{i_c} \in \mathbf{R}^{n_c}$ that has a single nonzero entry at the i_c th position. Then, consider the fine-grid function $\phi_{i_c}^{(H)}$ that has coefficient vector the i_c th column of P , i.e., equal to $P\mathbf{e}_{i_c}$. The set of functions $\{\phi_{i_c}^{(H)}\}$ forms the coarse basis of interest. The parameter H stands for characteristic size of the support of the coarse basis functions.

The above matrix-vector strong approximation property (1.1) admits the following finite element function form:

$$\|A\| a(u_h - u_H, u_h - u_H) \leq C_A \sum_{\mathbf{x}_i \in \mathcal{N}_h} f_i^2 = C_A \sum_{\mathbf{x}_i \in \mathcal{N}_h} \left(\int_{\Omega} f(\mathbf{x}) \varphi_i \, d\mathbf{x} \right)^2.$$

Using Cauchy-Schwarz inequality, we have

$$\sum_{\mathbf{x}_i \in \mathcal{N}_h} \left(\int_{\Omega} f(\mathbf{x}) \varphi_i \, d\mathbf{x} \right)^2 \leq \sum_{\mathbf{x}_i \in \mathcal{N}_h} \int_{\text{support}(\varphi_i)} f^2(\mathbf{x}) \, d\mathbf{x} \int_{\Omega} \varphi_i^2 \, d\mathbf{x}.$$

For a fairly general class of basis functions (including piecewise linears) on a quasiuniform mesh, we have

$$\int_{\Omega} \varphi_i^2 \, d\mathbf{x} \simeq |\text{support}(\varphi_i)| \simeq h^d.$$

Due to the bounded overlap of the supports of the finite element basis functions, we also have

$$\sum_{\mathbf{x}_i \in \mathcal{N}_h \text{ support } (\varphi_i)} \int f^2(\mathbf{x}) d\mathbf{x} \leq \kappa \|f\|_0^2.$$

Thus, we arrive at the energy error estimate of our main interest (using (1.2))

$$a(u_h - u_H, u_h - u_H) \leq C_A \kappa \frac{h^d}{\|A\|} \|f\|_0^2 \simeq C_A \kappa h^2 \|f\|_0^2.$$

In practice, we typically have $C_A = \mathcal{O}((\frac{H}{h})^2)$ with a constant in the \mathcal{O} symbol, independent of the two mesh sizes (h and H); see, e.g., Corollary 3.1 later on. Thus, we get the following final upscaling energy error estimate:

$$a(u_h - u_H, u_h - u_H) \leq c_A H^2 \|f\|_0^2. \quad (1.3)$$

In the remaining sections, we summarize a few AMG methods that provide *strong approximation property*.

2 Efficient window-based spectral AMG methods

In [FVZ05] (see also [Va08]), the following AMG method was proposed that exhibits strong approximation property. The original version tends to lead to relatively large coarse spaces so that the resulting two (and multi)-level methods have unacceptably high complexities. In the present section, we propose several approaches in the attempt to reduce the complexity of the original method.

Given an overlapping partition $\{w\}$ of the set of indices $i = 1, 2, \dots, n$, we extract the rows of a given $n \times n$ matrix A with indices from any given set (called window) w . The respective rectangular matrix is denoted by A_w . By proper reordering, A_w can be written as follows

$$A_w = [A_{ww}, A_{w,\chi}].$$

Here, A_{ww} is the principal submatrix of A (row and column indices from w) and $A_{w,\chi}$ is the submatrix of A with columns outside w (and row indices from w).

We are interested, for a proper nonnegative diagonal matrix D_w , in the normal matrices $A_w^T D_w A_w$. The diagonal matrices D_w provide a partition of unity, i.e., if I_w stands for extension by zero outside the set w , then $\sum_w I_w D_w I_w^T = I$. This property ensures that

$$\sum_w \mathbf{v}^T A_w^T D_w A_w \mathbf{v} = \sum_w \mathbf{v}^T A^T I_w D_w I_w^T A \mathbf{v} = \|A \mathbf{v}\|^2, \quad A_w = I_w^T A. \quad (2.1)$$

The method in question uses the symmetric semi-definite Schur complements S_w defined as follows:

$$\mathbf{v}_w^T S_w \mathbf{v}_w = \inf_{\mathbf{v}_\chi} \begin{bmatrix} \mathbf{v}_w \\ \mathbf{v}_\chi \end{bmatrix}^T A_w^T D_w A_w \begin{bmatrix} \mathbf{v}_w \\ \mathbf{v}_\chi \end{bmatrix}. \quad (2.2)$$

The original method utilizes the eigenvectors of the semidefinite Schur complements S_w ,

$$S_w \mathbf{p}_k = \lambda_k \mathbf{p}_k, \quad k = 1, \dots, n_w. \quad (2.3)$$

For efficiency reason, for a given tentative interpolation matrix \tilde{P} , we use in (2.3) instead the modified Schur complements

$$\mathbf{v}_w^T S_w \mathbf{v}_w = \inf_{\mathbf{v}_\chi \in I_\chi^T \text{Range}(\tilde{P})} \begin{bmatrix} \mathbf{v}_w \\ \mathbf{v}_\chi \end{bmatrix}^T A_w^T D_w A_w \begin{bmatrix} \mathbf{v}_w \\ \mathbf{v}_\chi \end{bmatrix}. \quad (2.4)$$

In what follows, we denote the exact window Schur complement with S_w^* .

We first form local interpolation matrices P_w by putting together the first $n_w^c \geq 1$ eigenvectors (in the lower part of the spectrum of S_w), i.e.,

$$P_w = [\mathbf{p}_1, \dots, \mathbf{p}_{n_w^c}]. \quad (2.5)$$

The corresponding eigenvalues (ordered in an increasing order) are such that $\lambda_k \leq \text{tol } \lambda_{\max}(S_w)$ for $k \leq n_w^c$ and

$$\lambda_k(S_w) > \text{tol } \|S_w\| = \text{tol } \lambda_{\max}(S_w) \quad \text{for } k > n_w^c. \quad (2.6)$$

Here, we have the freedom to choose the pre-selected tolerance “tol” (a number between zero and one) that may also vary with w .

The eigenvectors $\{\mathbf{p}_k\}_{k=1}^{n_w}$ are orthogonal and assumed normalized.

The global P is computed based on another partition of unity set of nonnegative $n_w \times n_w$ diagonal matrices $\{Q_w\}$ that satisfy

$$I = \sum_w I_w Q_w I_w^T.$$

Then P is defined as follows

$$P = \sum_w I_w Q_w [0, P_w, 0] = \sum_w I_w Q_w P_w (I_w^c)^T. \quad (2.7)$$

Here, I_w^c maps the local indices of the eigenvectors coming from the window w to their global indices expanding the result with zeros elsewhere. Thus we have defined a process that from a tentative \tilde{P} produces another one P . This can be iterated several times (by possibly changing the parameters such as $\{w\}$ and tol). In the next theorem, we formulate conditions ensuring that P admits a strong approximation property.

Theorem 2.1 *Consider the iterated window spectral AMG interpolation matrix P constructed on the basis of the modified window Schur complements using a \tilde{P} that satisfies the following estimate*

$$\sum_w \|D_w^{\frac{1}{2}} A_{w,\chi} (\mathbf{v}_\chi - I_\chi^T \tilde{P} \mathbf{w}_e)\|^2 \leq \mu \|A \mathbf{v}\|^2. \quad (2.8)$$

That is, \tilde{P} is such that for any \mathbf{v} , when restricted to a complementary set χ , there is a coarse vector \mathbf{w}_c (depending on \mathbf{v} and the set χ) such that for a fixed number $\mu > 0$ (2.8) holds. Then, if we choose $\text{tol} = \frac{1}{\delta} \leq 1$ in the two-level spectral decomposition defining the local P_w so that (see (2.6))

$$\|S_w\| \leq \delta \lambda_{m_w+1}(S_w),$$

and if we also assume the quasiuniformity of the windows, i.e., the estimate

$$\beta \|A\|^2 \leq \|S_w^*\|, \quad (2.9)$$

then, the following main strong approximation property holds for P

$$\|A\|^2 \|\mathbf{v} - P\mathbf{v}_c\|^2 \leq \eta \|A\mathbf{v}\|^2. \quad (2.10)$$

Here, $\eta = \frac{\delta}{\beta} (1 + \sqrt{\mu})^2$, where μ is from (2.8).

3 A new “window”-based spectral AMG method for finite element matrices

Here, we present a modified version of the method applied to finite element matrices A . The difference is in the eigenproblems that we use. Also, it utilizes a special partition of unity matrices. A main additional assumption is that the window sets are covered exactly by fine-grid elements and that we have access to the respective fine-grid element matrices so that we can assemble the semi-definite local matrices further denoted by Λ_w . Therefore, we have the identity $A = \sum_w I_w \Lambda_w I_w^T$.

We solve eigenproblems associated with the pair of matrices Λ_w and S_w , where S_w is the exact window-based Schur complement (as introduced before). The eigenproblems read (compare with (2.3)):

$$S_w \mathbf{p}_k = \lambda_k \Lambda_w \mathbf{p}_k, \quad k = 1, \dots, n_w, \quad (3.1)$$

where the eigenvalues are numbered in an increasing order and the eigenvectors are Λ_w -normalized.

Since, the matrices Λ_w can also be only semi-definite, to have real eigenvalues the nullspace of Λ_w should be contained in the nullspace of S_w , which is the case for finite elliptic matrices (Laplacian-like as well as elasticity).

It is clear that we can choose the eigenvectors \mathbf{p}_k be orthogonal to the nullspace of Λ_w (and Λ_w -orthogonal to each other). Let the columns of $P_w^{(0)}$ span the nullspace of Λ_w . Then, we have $\mathbf{p}_k^T P_w^{(0)} = 0$.

Based on a preselected tolerance $\text{tol} \in [0, 1)$, we choose n_w^c such that $\lambda_k > \text{tol} \lambda_{n_w}$ for $k > n_w^c$. The local interpolation matrices are defined similarly as before (cf., (2.5)), now augmented with the nullspace, i.e.,

$$P_w = \left[P_w^{(0)}, \mathbf{p}_1, \dots, \mathbf{p}_{n_w^c} \right]. \quad (3.2)$$

To define the global one, we use special diagonal matrices $\{Q_w\}_w$ with nonnegative entries that provide partition of unity, i.e., we have

$$I = \sum_w I_w Q_w I_w^T.$$

Each Q_w has entries on its diagonal $q_{w,i}$, $i \in w$, defined as follows:

$$q_{w,i} = \frac{\|\Lambda_w\|}{\sum_{w': i \in w'} \|\Lambda_{w'}\|}. \quad (3.3)$$

At the end we formulate our main result.

Theorem 3.1 *Let A be a given finite element s.p.d. matrix. Consider a given set of windows $\{w\}$ where each window w is exactly covered by fine-grid elements. Assume also that the local finite element matrices Λ_w corresponding to the sets w are available. The nullspace (if nonempty) of the local matrices Λ_w is assumed known (explicitly computed). That is, let the nullspace of Λ_w be represented by the range of an explicitly available local matrix $P_w^{(0)}$. Assume that this nullspace is contained in the nullspace of the window Schur complement S_w (defined in (2.2)). The global interpolation matrix P is defined as in (2.7) based on the local interpolation matrices (3.2) and the weights $q_{w,i}$ (entries of the diagonal partition of unity matrices Q_w) are defined in (3.3). Then, the following global strong approximation property holds*

$$\|\mathbf{v} - P\mathbf{v}_c\|_A^2 \leq \kappa \max_w \text{Cond}^+(\Lambda_w) \max_w \left(\frac{1}{\text{tol } \lambda_{\max}(\Lambda_w^+ S_w)} \right) \|A\mathbf{v}\|^2.$$

Above, $\kappa \geq 1$ depends on the overlap of $\{w\}$, $\text{tol} \in (0, 1]$ (in general depending on w) is the tolerance used to define the portion of the eigenvectors \mathbf{p}_k in the lower part of the spectrum computed in (3.1) used to define P_w , $\text{Cond}^+(\Lambda_w) = \frac{\|\Lambda_w\|}{\lambda_{\min}^+(\Lambda_w)}$ is the effective condition number of Λ_w computed in a subspace orthogonal to the nullspace of Λ_w . Finally, $\lambda_{\max}(\Lambda_w^+ S_w) = \max_k \lambda_k$ where λ_k are from (3.1).

Corollary 3.1 *For finite element s.p.d. matrices A coming from second order elliptic problems, the constructed finite element modification of the window-based spectral AMG method, the following strong approximation property holds*

$$\|\mathbf{v} - P\mathbf{v}_c\|_A^2 \leq C \kappa \left(\frac{H}{h} \right)^2 \max_w \frac{1}{\text{tol } \|A_{w_0, w_0}\|} \|A\mathbf{v}\|^2.$$

References

- [FVZ05] R. Falgout, P. S. Vassilevski, and L. T. Zikatanov, "On two-grid convergence estimates," Numerical Linear Algebra with Applications, **12**(2005), pp. 471–494.
- [Va08] Panayot S. Vassilevski, "MULTILEVEL BLOCK FACTORIZATION PRECONDITIONERS, Matrix-based Analysis and Algorithms for Solving Finite Element Equations," Springer, New York, 2008. 514 p.